Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081015\
Data File : BE090414.D
Acq On : 10 Aug 2015 22:36
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 11 03:56:53 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 06 12:28:08 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	132	0.00
2		1,4-Dioxane	0.687	0.766	-11.5	137	0.00
3		n-Nitrosodimethylamine	0.773	0.842	-8.9	144	0.00
4	S	2-Fluorophenol	0.750	1.175	-56.7#	221#	0.00
5	S	Phenol-d6	1.241	1.746	-40.7#	200#	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	138	0.00
7	S	Nitrobenzene-d5	0.331	0.375	-13.3	165#	0.00
8		Nitrobenzene	0.349	0.382	-9.5	169#	0.00
9		Naphthalene	1.104	1.127	-2.1	142	0.00
10		Hexachlorobutadiene	0.191	0.187	2.1	134	0.00
11		2-Methylnaphthalene	0.640	0.730	-14.1	165#	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	148	0.00
13	S	2,4,6-Tribromophenol	0.095	0.164	-72.6#	326#	0.00
14	S	2-Fluorobiphenyl	1.375	1.475	-7.3	156#	0.00
15		Acenaphthylene	8.098	8.880	-9.7	163#	0.00
16		Acenaphthene	1.259	1.269	-0.8	149	0.00
17		Fluorene	1.635	1.689	-3.3	151#	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	140	0.00
19		4-Bromophenyl-phenylether	0.194	0.207	-6.7	153#	0.00
20		Hexachlorobenzene	1.005	0.943	6.2	131	0.00
21		Pentachlorophenol	0.044	0.039	11.4	202#	0.00
22		Phenanthrene	1.251	1.269	-1.4	144	0.00
23		Anthracene	1.028	1.167	-13.5	164#	-0.02
24		Fluoranthene	1.231	1.410	-14.5	167#	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	141	0.00
26		Pyrene	1.084	1.252	-15.5	170#	0.00
27	S	Terphenyl-d14	0.737	0.865	-17.4	168#	0.00
28		Benzo(a)anthracene	1.108	1.202	-8.5	171#	0.00
29		Chrysene	1.339	1.373	-2.5	142	0.00
30		Bis(2-ethylhexyl)phthalate	0.311	0.458	-47.3#	290#	0.00
31		Indeno(1,2,3-cd)pyrene	0.942	1.054	-11.9	192#	0.00
32	I	Perylene-d12	1.000	1.000	0.0	152#	0.00
33		Benzo(b)fluoranthene	0.918	1.055	-14.9	186#	0.00
34		Benzo(k)fluoranthene	1.392	1.344	3.4	143	0.00
35	C	Benzo(a)pyrene	0.850	0.963	-13.3	183#	0.00
36		Dibenzo(a,h)anthracene	0.766	0.893	-16.6	205#	0.00
37		Benzo(g,h,i)perylene	0.923	1.030	-11.6	176#	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 0